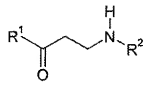
This listing of claims will replace all prior versions, and listings, of claims in the application:

LISTING OF CLAIMS:

1. **(Currently Amended)** A monoalkylaminoketone compound of the formula I



I

in which

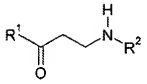
R^1 denotes a saturated, unsaturated or aromatic carbocyclic or heterocyclic radical which is unsubstituted or mono- or polysubstituted by R^3 and/or R^4, provided that R^1 is not 2,5-dimethyl-3-thienyl,

R^2 denotes alkyl having 1-20 C atoms,

R^3, R^4 each, independently of one another, denote H, alkyl or alkoxy having 1-20 C atoms, aryl, aryloxy or $COOR^2$, F, Br, OH, CN, NO_2, $N(R^2)_2$ or $NHCOR^2$,

or a salt ~~or solvate~~ thereof.

2. **(Withdrawn)** Process for the preparation of a monoalkylaminoketone compound of the formula I



I

in which

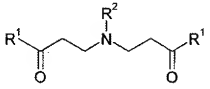
R^1 denotes a saturated, unsaturated or aromatic carbocyclic or heterocyclic radical which is unsubstituted or mono- or polysubstituted by R^3 and/or R^4, provided that R^1 is not 2,5-dimethyl-3-thienyl,

R^2 denotes alkyl having 1-20 C atoms,

R^3, R^4 each, independently of one another, denote H, alkyl or alkoxy having 1-20

C atoms, aryl, aryloxy or $COOR^2$, F, Br, OH, CN, NO_2, $N(R^2)_2$ or $NHCOR^2$,

by reacting a compound of the formula II



II

in which

R^1 and R^2 have the meaning indicated above, in the presence of an alkylamine of the formula R^2NH_2, in which R^2 has the meaning indicated above.

3. **(Withdrawn)** Process according to Claim 2, in which R^1 denotes phenyl or 2-thienyl.

4. **(Withdrawn)** Process according to Claim 2, in which R^2 denotes methyl, ethyl, n-propyl or isopropyl.

5. **(Withdrawn)** Process according to claim 2, wherein the pH for the conversion of the compounds of the formula II into the compounds of the formula I is adjusted to about pH 2-7.5 by addition of an alkylamine of the formula R^2NH_2.

6. **(Withdrawn)** Process according to claim 2, wherein the conversion of the compounds of the formula II into the compounds of the formula I is carried out at 0° - 200°C.

7. **(Withdrawn)** Process according to claim 2, wherein firstly the compound of the formula II is obtained by reaction of a mixture of a formaldehyde source with a corresponding alkylammonium salt and a ketone of the formula III

III

in which R^1 has the meaning indicated in Claim 1,
in the presence of a strong acid, and the compounds of the formula II obtained in this way are employed without further isolation for the preparation of the compounds of the

formula I.

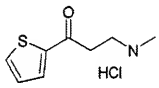
8. **(Withdrawn)** Process for the preparation of compounds of the formula I according to Claim 6, wherein the pH of the strongly acidic reaction mixture comprising the compounds of the formula II is increased to about pH 2-7.5, without further isolation of this compound, by addition of an alkylamine of the formula R^2NH_2, and the mixture is subsequently warmed.

9. **(Withdrawn)** Process for the preparation of compounds of the formula I according to Claim 7, wherein the reaction mixture comprising the compounds of the formula II is warmed to 0°C to 200°C after addition of a corresponding alkylamine.

10. **(Withdrawn)** Process according to claim 2 for the preparation of 3-methylamino-1-phenyl-1-propanone or 3-methylamino-1-(2-thienyl)-1-propanone.

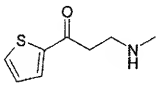
11. **(Withdrawn)** Process according to claim 2, wherein an acid-addition salt of the compound of the formula II is employed, and an acid-addition salt of the compound of the formula I is obtained.

12. **(Previously presented)** A compound of claim 1 which is of the formula Ia:



Ia

13. **(Currently Amended)** A compound of claim 1 which is of the formula Ib:



Ib

or a salt ~~or solvate~~ thereof.

14. (Canceled)

15. (Previously presented) A compound of claim 1, wherein R^1 denotes phenyl or 2-thienyl.

16. (Previously presented) A compound of claim 1, wherein R^2 denotes methyl, ethyl, n-propyl or isopropyl.

17. (Previously presented) A compound of claim 1, wherein R^1 is selected from: 2- or 3-furyl, 2- or 3-thienyl, 1-, 2- or 3-pyrrolyl, 1-, 2, 4- or 5-imidazolyl, 1-, 3-, 4- or 5-pyrazolyl, 2-, 4- or 5-oxazolyl, 3-, 4- or 5-isoxazolyl, 2-, 4- or 5-thiazolyl, 3-, 4- or 5-isothiazolyl, 2-, 3- or 4-pyridyl, 2-, 4-, 5- or 6-pyrimidinyl, furthermore preferably 1,2,3-triazol-1-, -4- or -5-yl, 1,2,4-triazol-1-, -3- or 5-yl, 1- or 5-tetrazolyl, 1,2,3-oxadiazol-4- or -5-yl, 1,2,4-oxadiazol-3- or -5-yl, 1,3,4-thiadiazol-2- or -5-yl, 1,2,4-thiadiazol-3- or -5-yl, 1,2,3-thiadiazol-4- or -5-yl, 3- or 4-pyridazinyl, pyrazinyl, 1-, 2-, 3-, 4-, 5-, 6- or 7-indolyl, 4- or 5-isoindolyl, 1-, 2-, 4- or 5-benzimidazolyl, 1-, 3-, 4-, 5-, 6- or 7-benzopyrazolyl, 2-, 4-, 5-, 6- or 7-benzoxazolyl, 3-, 4-, 5-, 6- or 7-benzisoxazolyl, 2-, 4-, 5-, 6- or 7-benzothiazolyl, 2-, 4-, 5-, 6- or 7-benzisothiazolyl, 4-, 5-, 6- or 7-benz-2,1,3-oxadiazolyl, 2-, 3-, 4-, 5-, 6-, 7- or 8-quinolyl, 1-, 3-, 4-, 5-, 6-, 7- or 8-isoquinolyl, 3-, 4-, 5-, 6-, 7- or 8-cinnolinyl, 2-, 4-, 5-, 6-, 7- or 8-quinazolinyl, 5- or 6-quinoxalinyl, 2-, 3-, 5-, 6-, 7- or 8-2H-benzo[1,4]oxazinyl, 1,3-benzodioxol-5-yl, 1,4-benzodioxan-6-yl, 2,1,3-benzothiadiazol-4- or -5-yl, 2,1,3-benzoxadiazol-5-yl, 2,3-dihydro-2-, -3-, -4- or -5-furyl, 2,5-dihydro-2-, -3-, -4- or 5-furyl, tetrahydro-2- or -3-furyl, 1,3-dioxolan-4-yl, tetrahydro-2- or -3-thienyl, 2,3-dihydro-1-, -2-, -3-, -4- or -5-pyrrolyl, 2,5-dihydro-1-, -2-, -3-, -4- or -5-pyrrolyl, 1-, 2- or 3-pyrrolidinyl, tetrahydro-1-, -2- or -4-imidazolyl, 2,3-dihydro-1-, -2-, -3-, -4- or -5-pyrazolyl, tetrahydro-1-, -3- or -4-pyrazolyl, 1,4-dihydro-1-, -2-, -3- or -4-pyridyl, 1,2,3,4-tetrahydro-1-, -2-, -3-, -4-, -5- or -6-pyridyl, 1-, 2-, 3- or 4-piperidinyl, 2-, 3- or 4-morpholinyl, tetrahydro-2-, -3- or -4-pyranyl, 1,4-dioxanyl, 1,3-dioxan-2-, -4- or -5-yl, hexahydro-1-, -3- or -4-pyridazinyl, hexahydro-1-, -2-, -4- or -5-pyrimidinyl, 1-, 2- or 3-piperazinyl, 1,2,3,4-tetrahydro-1-, -2-, -3-, -4-, -5-, -6-, -7- or -8-quinolyl, 1,2,3,4-tetrahydro-1-,-2-,-3-, -4-, -5-, -6-, -7- or -8-isoquinolyl, 2-, 3-, 5-, 6-, 7-

or 8- 3,4-dihydro-2H-benzo[1,4]oxazinyl, 2,3-methylenedioxyphenyl, 3,4-methylenedioxyphenyl, 2,3-ethylenedioxyphenyl, 3,4-ethylenedioxyphenyl, 3,4-(difluoromethylenedioxy)phenyl, 2,3-dihydrobenzofuran-5- or 6-yl, 2,3-(2-oxomethylenedioxy)phenyl, 3,4-dihydro-2H-1,5-benzodioxepin-6- or -7-yl, 2,3-dihydro-benzofuranyl or 2,3-dihydro-2-oxofuranyl,

each optionally substituted by R^3 and/or R^4.

18. (Previously presented) A compound of claim 1, wherein R^1 is selected from: phenyl; o-, m- or p-tolyl; o-, m- or p-hydroxyphenyl; o-, m- or p-methoxyphenyl; or, o-, m- or p-fluorophenyl.

19. (Previously presented) A compound of claim 1, wherein R^3 and R^4 are both H.